

~~orderable physicochemical property~~ corresponding to said each amino acid in said amino acid sequence of said target polypeptide or protein;

calculating one or more polypeptide eigenvalues and a corresponding polypeptide eigenvector associated with each of said one or more polypeptide eigenvalues by linear decomposition of an autocovariance matrix formed from a sequentially lagged data matrix of said polypeptide physicochemical data series;

ordering said one or more polypeptide eigenvalues and said corresponding polypeptide eigenvectors from largest to smallest;

selecting one or more of said polypeptide eigenvectors;

transforming said one or more of said polypeptide eigenvectors into an eigenvector template;

forming a graph of said eigenvector template, wherein said numerical values of ~~said physicochemical property~~ are graphed along the y-axis of said graph and ordered position in said eigenvector template is graphed along the x-axis of said graph;

partitioning said graph along said y-axis according to said ranges of said numerical values of ~~said physicochemical property~~ defining said peptide constituent groups, to form a plurality of y-axis ranges;

assigning one of said peptide constituents to each position in said peptide or peptide-like molecule by using said graph as a template to create a sequence of a mode-matched peptide, wherein at each ordered position in said eigenvector template along said x-axis of said graph, said one of said peptide constituents assigned to said ordered position has a numerical value of ~~said orderable physicochemical property~~ that is within said y-axis range of said ordered point;

determining a sequence of a retro-inverso peptide by inverting said sequence of a mode-matched peptide; and

synthesizing said retro-inverso peptide from said sequence, using D-amino acids.

2. A method for synthesizing a peptide or a peptide-like molecule based on matching a physicochemical mode of a peptide to the same physicochemical mode of a target polypeptide or protein, followed by synthesizing a retro-inverso version of said peptide comprised of D-amino acids, comprising the steps of:

assigning a numerical value of an orderable physicochemical property to each member of a set of peptide constituents a numerical value representative of the hydrophobic free energy of said each member of a set of peptide constituents, said set of peptide constituents including all the members of the set of naturally-occurring amino acids;

arranging said peptide constituents in order of said numerical values of said orderable physicochemical property;

partitioning said set of peptide constituents into a plurality of peptide constituent groups, whereby each of said peptide constituent groups contains at least one member of said set of peptide constituents, each peptide constituent group encompasses a range of said ordered numerical values, and each member of said set of peptide constituents belongs to only one peptide constituent group;

creating a polypeptide physicochemical data series by replacing each amino acid in an amino acid sequence with said numerical value of said orderable physicochemical property corresponding to said each amino acid in said amino acid sequence;

calculating one or more polypeptide eigenvalues and a corresponding polypeptide eigenvector associated with each of said one or more polypeptide eigenvalues by linear decomposition of an autocovariance matrix formed from a sequentially lagged data matrix of said polypeptide physicochemical data series;

ordering said one or more polypeptide eigenvalues and said corresponding polypeptide eigenvectors from largest to smallest;

selecting one or more of said polypeptide eigenvectors;

forming a vector, said vector being a sum of the products of each of said plurality of said polypeptide eigenvectors multiplied by the corresponding eigenvalue;

forming a graph of said vector, wherein said numerical values of said orderable physicochemical property are graphed along the y-axis of said graph, and ordered position in said eigenvector template is graphed along the x-axis of said graph;

partitioning said graph along said y-axis according to said range of said numerical values of said orderable physicochemical property defining said peptide constituent groups, to form a plurality of y-axis ranges; and

assigning one of said peptide constituents to each position in said peptide or peptide-like molecule by using said graph of said vector as a template, wherein at each ordered